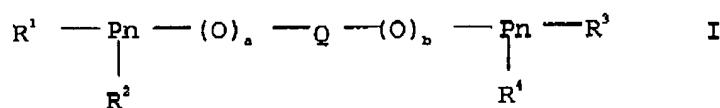


Application No. 10/527,635  
Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

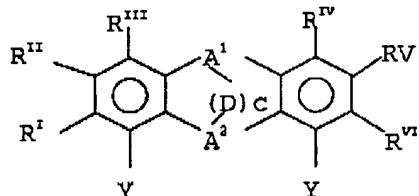
AMENDMENTS TO THE CLAIMS

1. (Original) A process for preparing dialdehydes and/or ethylenically unsaturated monoaldehydes by reacting at least one compound having at least two ethylenically unsaturated double bonds with carbon monoxide and hydrogen in the presence of a hydroformylation catalyst comprising at least one complex of a metal of transition group VIII with at least one ligand selected from among chelating pnicogen compounds of the formula I,



where

Q is a bridging group of the formula



where

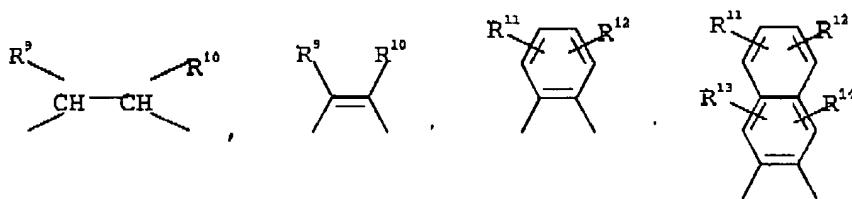
 $A^1$  and  $A^2$  are each, independently of one another, O, S,  $SiR^aR^b$ ,  $NR^c$  or  $CR^dR^e$ , where $R^a, R^b$  and  $R^e$  are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl or hetaryl,

Application No. 10/527,635  
Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

$R^d$  and  $R^e$  are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl or hetaryl or the group  $R^d$  together with a further group  $R^d$  or the group  $R^e$  together with a further group  $R^e$  form an intramolecular bridging group D,

D is a divalent bridging group selected from among the groups



where

$R^9$  and  $R^{10}$  are each, independently of one another, hydrogen, alkyl, cycloalkyl, aryl, halogen, trifluoromethyl, carboxyl, carboxylate or cyano or are joined to one another to form a C<sub>3</sub>-C<sub>4</sub>-alkylene bridge,

$R^{11}$ ,  $R^{12}$ ,  $R^{13}$  and  $R^{14}$  are each, independently of one another, hydrogen, alkyl, cycloalkyl, aryl, halogen, trifluoromethyl, COOH, carboxylate, cyano, alkoxy, SO<sub>3</sub>H, sulfonate, NE<sup>1</sup>E<sup>2</sup>, alkylene-NE<sup>1</sup>E<sup>2</sup>E<sup>3+</sup>X<sup>-</sup>, acyl or nitro,

c 0 or 1,

Y is a chemical bond,

$R^I$ ,  $R^{II}$ ,  $R^{III}$ ,  $R^{IV}$ ,  $R^V$  and  $R^{VI}$  are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl, hetaryl, COOR<sup>f</sup>, COO<sup>-</sup>M<sup>+</sup>, SO<sub>3</sub>R<sup>f</sup>, SO<sub>3</sub><sup>-</sup>M<sup>+</sup>,

Application No. 10/527,635  
Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

NE<sup>1</sup>E<sup>2</sup>, NE<sup>1</sup>E<sup>2</sup>E<sup>3+</sup>X<sup>-</sup>, alkylene-NE<sup>1</sup>E<sup>2</sup>E<sup>3+</sup>X<sup>-</sup>, OR<sup>f</sup>, SR<sup>f</sup>, (CHR<sup>g</sup>CH<sub>2</sub>O)<sub>x</sub>R<sup>f</sup>,  
(CH<sub>2</sub>N(E<sup>1</sup>))<sub>x</sub>R<sup>f</sup>, (CH<sub>2</sub>CH<sub>2</sub>N(E<sup>1</sup>))<sub>x</sub>R<sup>f</sup>, halogen, trifluoromethyl, nitro, acyl or cyano,

where

R<sup>f</sup>, E<sup>1</sup>, E<sup>2</sup> and E<sup>3</sup> are identical or different radicals selected from among hydrogen, alkyl,  
cycloalkyl and aryl,

R<sup>g</sup> is hydrogen, methyl or ethyl,

M<sup>+</sup> is a cation,

X<sup>-</sup> is an anion, and

x is an integer from 1 to 120,

or

two adjacent radicals selected from among R<sup>I</sup>, R<sup>II</sup>, R<sup>III</sup>, R<sup>IV</sup>, R<sup>V</sup> and R<sup>VI</sup> together with  
two adjacent carbon atoms of the benzene ring to which they are bound for a fused  
ring system having 1, 2 or 3 further rings,

a and b are each, independently of one another, 0 or 1,

Pn is a pnicogen atom selected from among the elements phosphorus, arsenic and  
antimony,

and

Application No. 10/527,635  
Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

$R^1, R^2, R^3, R^4$  are each, independently of one another, hetaryl, hetaryloxy, alkyl, alkoxy, aryl, aryloxy, cycloalkyl, cycloalkoxy, heterocycloalkyl, heterocycloalkoxy or an  $NE^1E^2$  group, with the proviso that  $R^1$  and  $R^3$  are pyrrole groups bound via the nitrogen atom to the pnicogen atom  $Pn$

or  $R^1$  together with  $R^2$  and/or  $R^3$  together with  $R^4$  form a divalent group E of the formula

Py-I-W

where

Py is a pyrrole group which is bound via the pyrrole nitrogen atom to the pnicogen atom  $Pn$ ,

I is a chemical bond or O, S,  $SiR^aR^b$ ,  $NR^c$ , substituted or unsubstituted  $C_1-C_{10}$ -alkylene or  $CR^hR^i$ ,

W is cycloalkyl, cycloalkoxy, aryl, aryloxy, hetaryl or hetaryloxy,

and

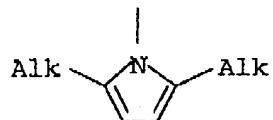
$R^h$  and  $R^i$  are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl or hetaryl,

or  $R^1$  together with  $R^2$  and/or  $R^3$  together with  $R^4$  form a bispyrrole group of the formula

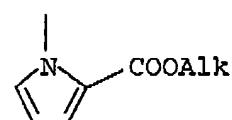
Py-I-Py

bound via the nitrogen atoms to the pnicogen atom Pn.

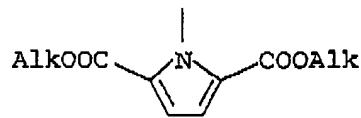
2. (Original) A process as claimed in claim 1, wherein at least one ligand of the formula I, in which the radicals R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> are selected independently from among groups of the formulae I.a to I.k



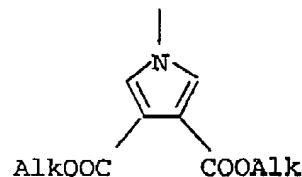
(I.a)



(I.b)



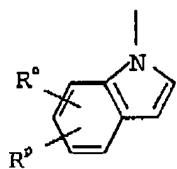
(I.c)



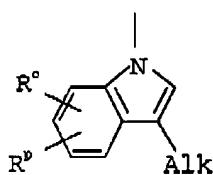
(I.d)

Application No. 10/527,635  
 Reply to Office Action of January 27, 2006

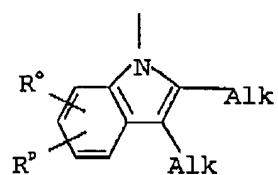
Docket No.: 13111-00007-US



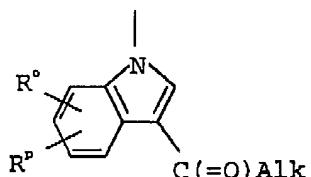
(I.e)



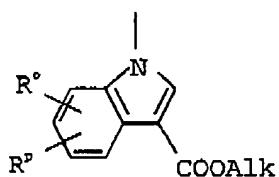
(I.f)



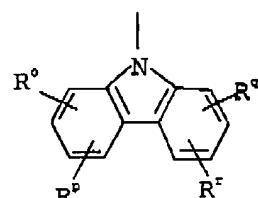
(I.g)



(I.h)



(I.i)



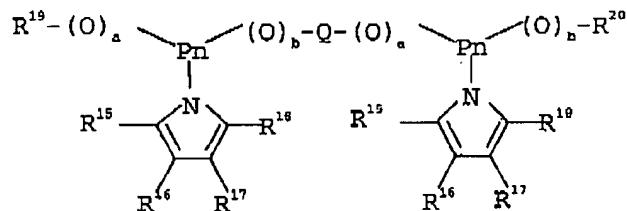
(I.k)

where

Alk is a C<sub>1</sub>-C<sub>4</sub>-alkyl group and

R°, R°, R° and R° are each, independently of one another, hydrogen, C<sub>1</sub>-C<sub>4</sub>-alkyl, C<sub>1</sub>-C<sub>4</sub>-alkoxy, acyl, halogen, trifluoromethyl, C<sub>1</sub>-C<sub>4</sub>-alkoxycarbonyl or carboxyl, is used.

3. (Currently amended) A process as claimed in claim 2, wherein at least one ligand of the formula I, in which the radicals R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> are each, independently of one another, a 3-alkylindolyl group, preferably a 3-methylindolyl group, is used.
4. (Previously presented) A process as claimed in claim 1, wherein the chelating pnicogen compound of the formula I is selected from among chelating pnicogen compounds of the formula II,



(II)

where

R<sup>15</sup>, R<sup>16</sup>, R<sup>17</sup> and R<sup>18</sup> are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl, hetaryl, W'COOR<sup>k</sup>, W'COO'M<sup>+</sup>, W'(SO<sub>3</sub>)R<sup>k</sup>, W'(SO<sub>3</sub>)M<sup>+</sup>, W'PO<sub>3</sub>(R<sup>k</sup>)(R<sup>l</sup>), W'(PO<sub>3</sub>)<sup>2-</sup>(M<sup>+</sup>)<sub>2</sub>, W'NE<sup>4</sup>E<sup>5</sup>, W'(NE<sup>4</sup>E<sup>5</sup>E<sup>6</sup>)<sup>+</sup>X<sup>-</sup>, W'OR<sup>k</sup>, W'SR<sup>k</sup>, (CHR<sup>l</sup>CH<sub>2</sub>O)<sub>y</sub>R<sup>k</sup>, (CH<sub>2</sub>NE<sup>4</sup>)<sub>y</sub>R<sup>k</sup>, (CH<sub>2</sub>CH<sub>2</sub>NE<sup>4</sup>)<sub>y</sub>R<sup>k</sup>, halogen, trifluoromethyl, nitro, acyl or cyano,

where

W' is a single bond, a heteroatom or a divalent bridging group having from 1 to 20 bridge atoms,

Application No. 10/527,635  
Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

$R^k$ ,  $E^4$ ,  $E^5$ ,  $E^6$  are identical or different radicals selected from among hydrogen, alkyl, cycloalkyl and aryl,

$R^1$  is hydrogen, methyl or ethyl,

$M^+$  is a cation equivalent,

$X^-$  is an anion equivalent and

$y$  is an integer from 1 to 240,

where two adjacent radicals  $R^{15}$ ,  $R^{16}$ ,  $R^{17}$  and  $R^{18}$  together with the carbon atoms of the pyrrole ring to which they are bound may also form a fused ring system having 1, 2 or 3 further rings,

with the proviso that at least one of the radicals  $R^{15}$ ,  $R^{16}$ ,  $R^{17}$  and  $R^{18}$  is not hydrogen and  $R^{19}$  and  $R^{20}$  are not joined to one another,

$R^{19}$  and  $R^{20}$  are each, independently of one another, cycloalkyl, heterocycloalkyl, aryl or hetaryl, or  $R^{19}$  together with  $R^{15}$  or  $R^{16}$  and/or  $R^{19}$  together with  $R^{17}$  or  $R^{18}$  form a divalent group

-I-W-

where

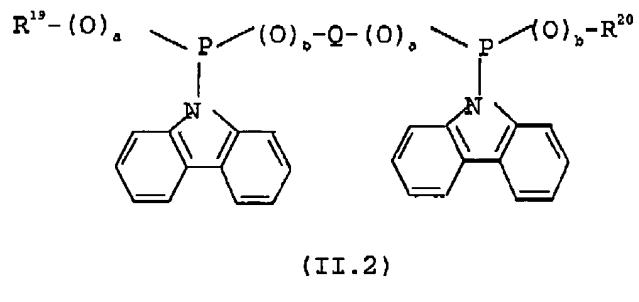
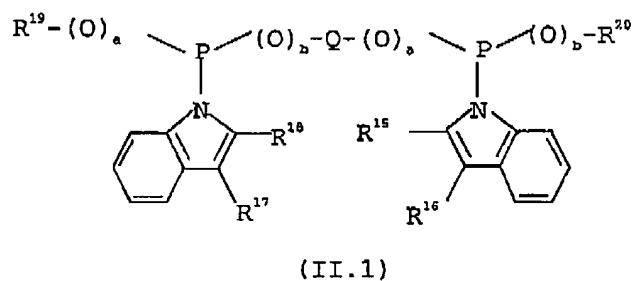
$I$  is a chemical bond or O, S,  $SiR^aR^b$ ,  $NR^c$  or substituted or unsubstituted  $C_1-C_{10}$ -alkylene, preferably  $CR^hR_i$ , where  $R^a$ ,  $R^b$ ,  $R^c$ ,  $R^h$  and  $R^i$  are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl or hetaryl and

Application No. 10/527,635  
Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

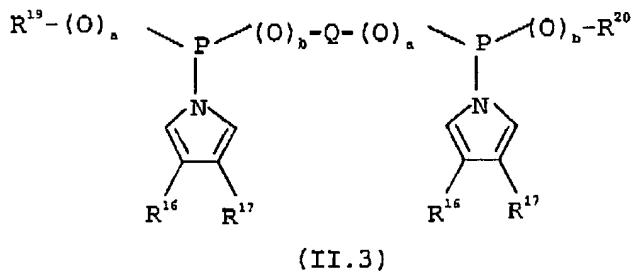
**W** is cycloalkyl, cycloalkoxy, aryl, aryloxy, hetaryl or hetaryloxy.

5. (Currently amended) A process as claimed in claim 1, wherein the chelating pnicogen compound of the formula I is a chelating pnicogen compound of the formulae II.1 to II.3,



Application No. 10/527,635  
 Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US



where

R<sup>15</sup>, R<sup>16</sup>, R<sup>17</sup>, R<sup>18</sup>, Q, a and b are as defined in claim 4,  
Q, a and b are as defined in claim 1,

R<sup>15</sup>, R<sup>16</sup>, R<sup>17</sup> and R<sup>18</sup> are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl, hetaryl, W'COOR<sup>k</sup>, W'COOM<sup>+</sup>, W'(SO<sub>3</sub>)R<sup>k</sup>, W'(SO<sub>3</sub>)M<sup>+</sup>, W'PO<sub>3</sub>(R<sup>k</sup>)(R<sup>l</sup>), W'(PO<sub>3</sub>)<sup>2-</sup>(M<sup>+</sup>)<sub>2</sub>, W'NE<sup>4</sup>E<sup>5</sup>, W'(NE<sup>4</sup>E<sup>5</sup>)<sup>2-</sup>X, W'OR<sup>k</sup>, W'SR<sup>k</sup>, (CHR<sup>1</sup>CH<sub>2</sub>O)<sub>y</sub>R<sup>k</sup>, (CH<sub>2</sub>NE<sup>4</sup>)<sub>y</sub>R<sup>k</sup>, (CH<sub>2</sub>CH<sub>2</sub>NE<sup>4</sup>)<sub>y</sub>R<sup>k</sup>, halogen, trifluoromethyl, nitro, acyl or cyano,

wherein

W' is a single bond, a heteroatom or a divalent bridging group having from 1 to 20 bridge atoms.

R<sup>k</sup>, E<sup>4</sup>, E<sup>5</sup>, E<sup>6</sup> are identical or different radicals selected from among hydrogen, alkyl, cycloalkyl and aryl,

Application No. 10/527,635  
 Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

R<sup>1</sup> is hydrogen, methyl or ethyl.

M<sup>+</sup> is a cation equivalent.

X is an anion equivalent and

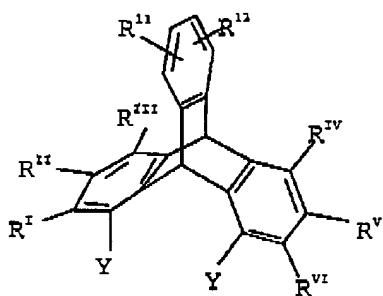
y is an integer from 1 to 240,

where two adjacent radicals R<sup>15</sup>, R<sup>16</sup>, R<sup>17</sup> and R<sup>18</sup> together with the carbon atoms of the pyrrole ring to which they are bound may also form a fused ring system having 1, 2 or 3 further rings.

where at least one of the radicals R<sup>16</sup> and R<sup>17</sup> in the formula II.3 is not hydrogen,

R<sup>19</sup> and R<sup>20</sup> are each, independently of one another, cycloalkyl, heterocycloalkyl, aryl or hetaryl.

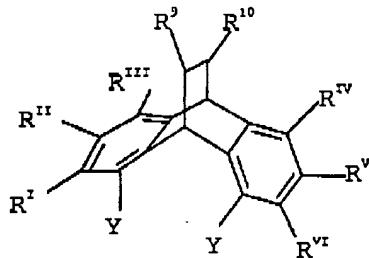
6. (Previously presented) A process as claimed in claim 1, wherein the bridging group Q is a triptycenediyl group of the formula



Application No. 10/527,635  
 Reply to Office Action of January 27, 2006

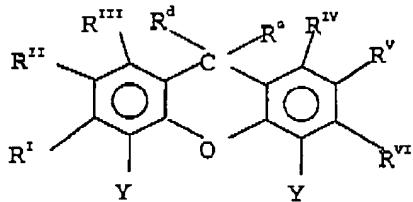
Docket No.: 13111-00007-US

or the formula



where R<sup>I</sup>, R<sup>II</sup>, R<sup>III</sup>, R<sup>IV</sup>, R<sup>V</sup> and R<sup>VI</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup> and R<sup>12</sup> are as defined in claim 1.

7. (Previously presented) A process as claimed in claim 1, wherein the bridging group Q is a xanthenediyl group of the formula



where R<sup>I</sup>, R<sup>II</sup>, R<sup>III</sup>, R<sup>IV</sup>, R<sup>V</sup> and R<sup>VI</sup> and Y are as defined in claim 1 and R<sup>d</sup> and R<sup>e</sup> are each, independently of one another, hydrogen, alkyl, cycloalkyl, heterocycloalkyl, aryl or hetaryl.

8. (Previously presented) A process as claimed in claim 1, wherein a molar ratio of ligand to metal of transition group VIII of from 1:1 to 1000:1 is set in the reaction mixture.

Application No. 10/527,635  
Reply to Office Action of January 27, 2006

Docket No.: 13111-00007-US

9. (Previously presented) A process as claimed in claim 1, wherein the reaction is carried out at from 40 to 80°C.
10. (Previously presented) A process as claimed in claim 1, wherein the compound having at least two ethylenically unsaturated double bonds which is used is a  $\alpha,\omega$ -diolefin.
11. (Previously presented) A process as claimed in claim 1, wherein
  - (i) a compound having a least two ethylenically unsaturated double bonds is subjected to the hydroformylation reaction in a reaction zone,
  - (ii) an output is taken from the reaction zone and is fractionated to give a fraction enriched in unsaturated monoaldehydes and a fraction depleted in unsaturated monoaldehydes, and
  - (iii) the fraction depleted in unsaturated monoaldehydes is recirculated, optionally after work up, to the reaction zone.
12. (New) A process as claimed in claim 2, wherein at least one ligand of the formula I, in which the radicals R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> are each, independently of one another, a 3-methylindolyl group.